

WEST Search History

DATE: Thursday, March 20, 2003

Set Name Query
side by side

Hit Count Set Name
result set

*DB=USPT,PGPB,JPAB,EPAB,DWPI,TDBD; THES=ASSIGNEE;
PLUR=YES; OP=AND*

L3 L2 and alpha

204 L3

L2 L1 and (alcohol\$3 or alkanol\$3)

660 L2

L1 (carbonyl\$3 or aldehyde\$3 or ketone\$3) near15 (electrochem\$6 or
electroly\$6 or electro\$2oxid\$8)

1363 L1

END OF SEARCH HISTORY

=> d his

(FILE 'HOME' ENTERED AT 11:38:09 ON 20 MAR 2003)

FILE 'CAPLUS' ENTERED AT 11:38:26 ON 20 MAR 2003

L1	16294 S (CARBONYL# OR ALDEHYDE# OR KETONE#) (15A) ELECTRO?
L2	1053 S L1 AND (ALCOHOL# OR ALKANOL#)
L3	172 S L2 AND ALPHA
L4	22 S L2 AND (ALPHA (2A) HYDROXY?)

=>

=> d

L3 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2003 ACS
RN 1320-67-8 REGISTRY
CN Propanol, 1(or 2)-methoxy- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Propanol, methoxy- (7CI, 8CI)
OTHER NAMES:
CN Arcosolv PM
CN Dowanol PM
CN Kuraray PGM
CN **Methoxypropanol**
CN Methyl Propasol
CN PM
CN PM (solvent)
CN Polysolv MPM
CN Propylene glycol methyl ether
CN Propylene glycol monomethyl ether
CN Solvenon PM
DR 170516-38-8, 1337-64-0, 143749-19-3, 89024-56-6, 30523-02-5
MF C4 H10 O2
CI IDS, COM
LC STN Files: BIOBUSINESS, BIOSIS, CA, CAOLD, CAPLUS, CASREACT, CEN,
CHEMCATS, CHEMLIST, CIN, CSCHEM, CSNB, DETHERM*, EMBASE, IFICDB,
IFIPAT,
IFIUDB, MEDLINE, MSDS-OHS, NIOSHTIC, PDLCOM*, PIRA, PROMT, RTECS*,
TOXCENTER, ULIDAT, USPAT2, USPATFULL, VTB
(*File contains numerically searchable property data)
Other Sources: DSL**, EINECS**
(**Enter CHEMLIST File for up-to-date regulatory information)

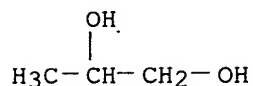
CM 1

CRN 67-56-1
CMF C H4 O

H₃C-OH

CM 2

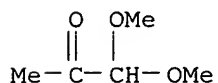
CRN 57-55-6
CMF C3 H8 O2



1208 REFERENCES IN FILE CA (1962 TO DATE)
36 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
1212 REFERENCES IN FILE CAPLUS (1962 TO DATE)
1 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

=> d

L2 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2003 ACS
RN 6342-56-9 REGISTRY
CN 2-Propanone, 1,1-dimethoxy- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Pyruvaldehyde, 1-(dimethyl acetal) (6CI, 8CI)
OTHER NAMES:
CN 1,1-Dimethoxy-2-propanone
CN 1,1-Dimethoxyacetone
CN Dimethoxymethyl methyl ketone
CN Methylglyoxal dimethyl acetal
CN Pyruvaldehyde dimethyl acetal
CN Pyruvic aldehyde dimethyl acetal
FS 3D CONCORD
MF C5 H10 O3
LC STN Files: BEILSTEIN*, BIOBUSINESS, BIOSIS, CA, CAOLD, CAPLUS,
CASREACT,
CHEMCATS, CHEMINFORMRX, CHEMLIST, CSCHEM, IFICDB, IFIPAT, IFIUDB,
SPECINFO, SYNTHLINE, TOXCENTER, USPAT2, USPATFULL
(*File contains numerically searchable property data)
Other Sources: EINECS**, NDSL**, TSCA**
(**Enter CHEMLIST File for up-to-date regulatory information)



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

337 REFERENCES IN FILE CA (1962 TO DATE)
338 REFERENCES IN FILE CAPLUS (1962 TO DATE)
2 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

=>